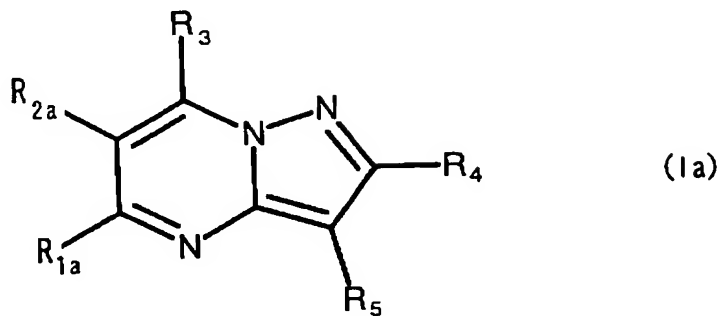


What is Claimed is:

1. A compound represented by the formula:



or a prodrug, pharmaceutically acceptable salt or solvate thereof,

- 5 wherein, in the formula (Ia), R_{1a} , R_{2a} , R_3 - R_5 represent, each independently, hydrogen, halogen, lower alkyl that may be substituted, lower alkenyl that may be substituted, lower alkynyl that may be substituted, cycloalkyl that may be substituted, cycloalkenyl that may be substituted, cycloalkynyl that may be substituted, aryl that may be substituted, heterocyclic group that may be substituted, hydroxy, alkoxy that may be substituted, aryloxy that may be substituted, heterocyclic oxy that may be substituted, acyl that may be substituted, monosubstituted carbonyloxy that may be substituted, carbamoyl that may be substituted, diazo, amidino that may be substituted, azido, nitroso, nitro, amino that may be substituted, imino that may be substituted, cyano, mercapto, monosubstituted thio that may be substituted, monosubstituted thioxy that may be substituted, monosubstituted sulfinyl that may be substituted, monosubstituted sulfonyl that may be substituted,
- 10
- 15
- 20

25 sulfo, or trisubstituted silyl, and any combinations of
R_{1a}, R_{2a}, R₃-R₅ may together form a ring structure;
provided that the following (i)-(x) are excluded:

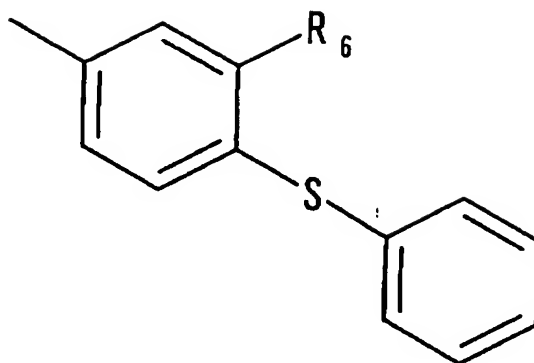
(i) a compound, wherein R_{1a} is hydrogen,
OH, lower alkyl, cycloalkyl having a carbon number of
3-8, halogenated lower alkyl, or phenyl;

30 R_{2a} is hydrogen, lower alkoxy carbonyl,
lower alkoxy, halogen, lower alkyl, cycloalkyl having a
carbon number of 3-8, lower alkoxy carbonyl lower alkyl,
carboxyl, carboxy lower alkyl, -CONHR₆ (R₆: hydrogen;
phenyl that may have a halogen atom, or lower alkyl),
35 cyano; phenyl that may have a substituent selected from
the group consisting of a hydroxyl group, halogen atom,
lower alkyl group, lower alkoxy and phenylthio group;
phenyl lower alkyl group that may have a substituent
selected from the group consisting of hydroxyl group
40 and lower alkoxy group on the phenyl ring; lower
alkanoyloxy lower alkyl; benzoyl group; lower alkanoyl
group that may have halogen atom; or hydroxy lower
alkyl group that may have a substituent selected from
the group consisting of a phenyl group and halogen
45 atom;

R₃ is hydrogen, or OH;

R₄ is hydrogen, lower alkyl, lower alkoxy
lower alkyl, or halogenated lower alkyl;

R₅ is



50

and

R_6 is hydrogen, lower alkyl, or lower alkoxy;

(ii) a compound, wherein R_{1a} and R_{2a} are,
55 each independently, hydrogen, halogen, CN, alkyl, alkenyl, alkynyl, cycloalkyl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, amino, alkylamino, or (substituted) phenyl; and

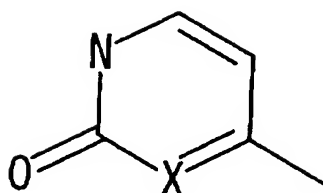
R_3 is (substituted) aryl, or
60 (substituted) heteroaryl;

(iii) a compound, wherein R_{1a} is hydrogen, (substituted) lower alkyl, cycloalkyl, thienyl, furyl, lower alkenyl, or (substituted) phenyl;

R_{2a} is hydrogen or lower alkyl; and
65 R_3 is amino that may be substituted;

(iv) a compound, wherein R_{1a} is hydrogen, alkyl, OH, O-alkyl, halo, amino, or nitro;

R_{2a} is



70 wherein X is CH or N, and the nitrogen atom on the R_{2a} ring may be substituted; and

R₃ and R₅ are, each independently, hydrogen, alkyl, alkenyl, alkynyl, aryl, halo, OH, or heterocyclyl;

75 (v) a compound, wherein R_{1a} is hydrogen, alkyl, alkoxy, OH, halo, NO₂, or NH₂;

R_{2a} is hydrogen, (substituted) alkyl, cycloalkyl, alkoxy, (substituted) alkenyl, (substituted) alkynyl, (substituted) aryl, 80 (substituted) heterocyclyl, alkoxy-NRR, NO₂, OH, NH₂, or (substituted) heteroaryl;

R₃ and R₄ are, each independently, hydrogen, alkyl, aryl, cycloalkyl, OH, halo, amino, or nitro; and

85 R₅ is hydrogen, (substituted) alkyl, cycloalkyl, aryl, (substituted) heterocyclyl, halo, OH, or (substituted) heteroaryl;

(vi) a compound, wherein R_{2a} is substituted acetyl, or heterocyclic-substituted lower 90 alkylene or lower alkenylene; and

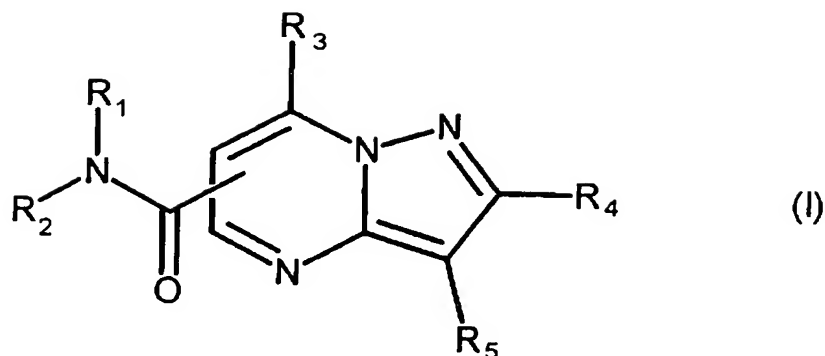
R₃ is phenyl that may be substituted;

(vii) a compound, wherein R_{1a} and R_{2a} are each independently, hydrogen, halogen, (substituted) alkyl, (substituted) alkenyl, (substituted) aryl,

95 (substituted) aralkyl, (substituted) heterocyclic
group, or together form an alkylene group; and
R₃ is amino that may be substituted;
(viii) a compound, wherein R_{1a} is
hydrogen, alkyl, cycloalkyl, alkoxy, (alkyl)amino,
100 aryl, or heteroaryl;
R_{2a} is hydrogen, alkyl, halogen, cyano,
hydroxy, or alkoxy;
R₃ is amino that may be substituted, or
alkoxy that may be substituted; and
105 R₅ is aryl;
(ix) R_{1a} is lower alkyl that is
substituted with a substituent selected from the group
consisting of carboxy, lower alkoxy-carboxy, and
substituted carbamoyl;
110 R_{2a} is hydrogen;
R₃ is phenylcarbonylamino, wherein said
phenyl group may be substituted; and
R₄ and R₅ are hydrogen;
(x) (2, 5-dimethyl-pyrazolo-[1, 5-a]-
115 pyrimidine-7-yl) succinic acid;
wherein the undefined substituents in
the compounds (i)-(x) represent any substituents.

2. The compound of claim 1, wherein either
one of R_{1a} and R_{2a} is hydrogen, and the other one is
carbamoyl that may be substituted.

3. The compound of claim 1, represented by
the formula:



or a prodrug, pharmaceutically acceptable salt or
5 solvate thereof,

wherein, in the formula (I),

R₁ is hydrogen, lower alkyl, amino that
may be substituted, or aryl lower alkyl that may be
substituted; and

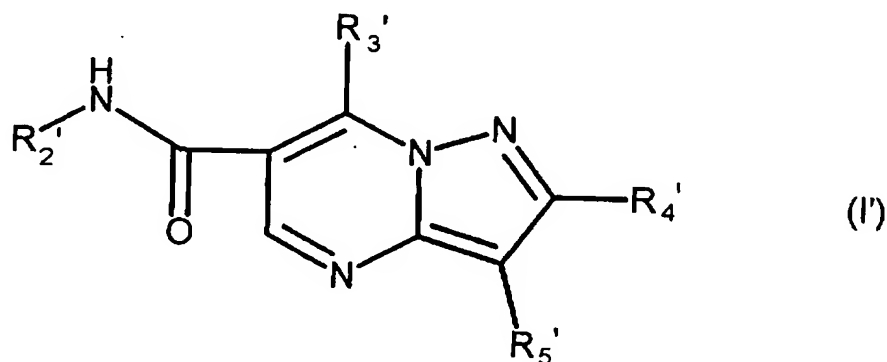
10 R₂ is hydrogen, lower alkyl that may be
substituted, cycloalkyl that may be substituted,
cycloalkyl lower alkyl that may be substituted, lower
alkoxy that may be substituted, aryl that may be
substituted, aryl lower alkyl that may be substituted,
15 aryloxy lower alkyl that may be substituted, lower
alkylsulfonyl that may be substituted, arylsulfonyl
that may be substituted, heteroaryl lower alkyl that
may be substituted, heterocyclic group lower alkyl that
may be substituted, or amino that may be substituted;
20 or

R₁ and R₂ together with the adjacent N
atom may form a heterocycle that may be substituted;

R₃ is hydrogen, hydroxy, lower alkoxy,
halogen, or amino that may be substituted;

25 R₄ is hydrogen, lower alkyl, or aryl that
may be substituted; and

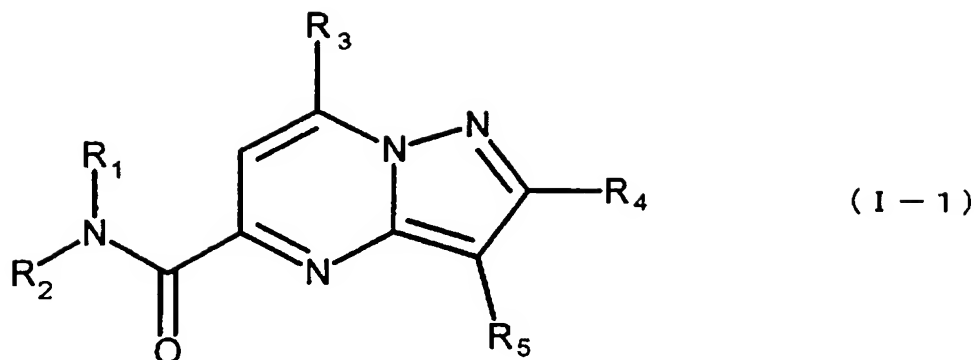
R₅ is hydroxy, lower alkyl that may be substituted, aryl that may be substituted, aryl lower alkyl that may be substituted, cycloalkyl lower alkyl that may be substituted, aryl lower alkenyl that may be substituted, cycloalkyl lower alkenyl that may be substituted, aryl lower alkynyl that may be substituted, cycloalkyl lower alkynyl that may be substituted, aryl carbonyl that may be substituted, aryl lower alkyl carbonyl that may be substituted, heterocyclic group that may be substituted, halogen, CHO, amino that may be substituted, or imino that may be substituted; provided that a compound represented by the following formula is excluded:



40

wherein, in the formula (I'),
R₂' is hydrogen, phenyl that may be substituted with lower alkyl or halogen; R₃' is hydrogen or hydroxy; R₄' is hydrogen or lower alkyl; and R₅' is phenyl having phenylthio group that may further be substituted with lower alkyl or lower alkoxy.

4. The compound of claim 3 represented by the formula:



5 or a prodrug, pharmaceutically acceptable salt or solvate thereof,

wherein, in the formula (I-1), each substituent is defined above.

5. The compound of claim 3 or 4, or a prodrug, pharmaceutically acceptable salt or solvate thereof, wherein R_1 is hydrogen; and R_2 is aryl that may be substituted.

6. The compound of claim 3 or 4, or a prodrug, pharmaceutically acceptable salt or solvate thereof, wherein R_3 is hydrogen, or amino that may be substituted.

7. The compound of claim 3 or 4, or a prodrug, pharmaceutically acceptable salt or solvate thereof, wherein R_4 is hydrogen.

8. The compound of claim 3 or 4, or a prodrug, pharmaceutically acceptable salt or solvate thereof, wherein R_5 is aryl that may be substituted.

9. The compound of claim 3 or 4, or a prodrug, pharmaceutically acceptable salt or solvate thereof, wherein R_1 is hydrogen; R_2 is phenyl that may

be substituted; R_3 is hydrogen, or amino that may be
5 substituted; R_4 is hydrogen; and R_5 is phenyl that may be substituted.

10. The compound of claim 9, or a prodrug, pharmaceutically acceptable salt or solvate thereof, wherein the substituent on the phenyl in R_2 that may be substituted is one or more selected from the group
5 consisting of heterocyclic group that may be substituted, lower alkyl carbonyl, cycloalkyl, lower alkyl, amino that may be substituted, halogen, halogenated lower alkyl, lower alkoxy, carboxy lower alkyloxy, heterocyclic group lower alkyloxy, amino
10 lower alkyl, hydroxy, cyano, carbamoyl-heterocyclic group-oxy, cyano lower alkyl, and phenyl.

11. The compound of claim 10, or a prodrug, pharmaceutically acceptable salt or solvate thereof, wherein R_2 is heterocyclic group phenyl that may be substituted.

12. The compound of claim 10, or a prodrug, pharmaceutically acceptable salt or solvate thereof, wherein R_2 is piperazino phenyl that may be substituted, piperizino phenyl that may be substituted, or
5 pyrrolidino phenyl that may be substituted.

13. The compound of claim 9, or a prodrug, pharmaceutically acceptable salt or solvate thereof, wherein the substituent on the phenyl in R_5 that may be substituted is one or more selected from the group
5 consisting of halogen, halogenated lower alkyl, aryl lower alkyloxy, lower alkyl, lower alkoxy, hydroxy, lower alkylthio, phenyl, phenyloxy, phenyl lower alkyl, phenyl lower alkylamino, phenyl lower alkylthio, phenyl

lower alkenyl, phenyl carbamoyl, amino, cycloalkyl
10 lower alkyloxy, and heteroaryl lower alkyloxy.

14. A pharmaceutical composition, comprising
the compound of any one of claims 1-13.

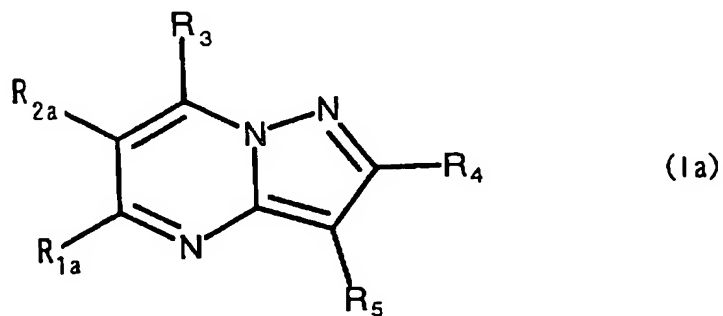
15. A NAD(P)H oxydase inhibitor, comprising
the compound of any one of claims 1-13.

16. A prophylactic or therapeutic agent for
NAD(P)H-related diseases, comprising the compound of
any one of claims 1-13.

17. The prophylactic or therapeutic agent of
claim 16, wherein said disease is selected from the
group consisting of inflammation, pulmonary circulation
disorders, ischemic heart disease, cerebral circulation
5 disorders, arteriosclerosis, diabetic complications,
hypertension, and proliferative disorders.

18. The prophylactic or therapeutic agent of
claim 16, wherein said disease is brain infarction or
diabetic retinal disorder.

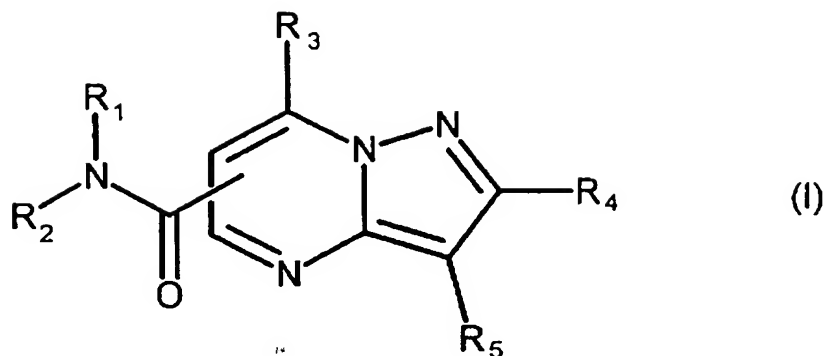
19. A NAD(P)H oxydase inhibitor, comprising
a compound represented by the formula (Ia):



or a prodrug, pharmaceutically acceptable salt or
5 solvate thereof,

wherein, in the formula, R_{1a} , R_{2a} , R_3-R_5
represent, each independently, hydrogen, halogen, lower
alkyl that may be substituted, lower alkenyl that may
be substituted, lower alkynyl that may be substituted,
10 cycloalkyl that may be substituted, cycloalkenyl that
may be substituted, cycloalkynyl that may be
substituted, aryl that may be substituted, heterocyclic
group that may be substituted, hydroxy, alkoxy that may
be substituted, aryloxy that may be substituted,
15 heterocyclic oxy that may be substituted, acyl that may
be substituted, monosubstituted carbonyloxy that may be
substituted, carbamoyl that may be substituted, diazo,
amidino that may be substituted, azido, nitroso, nitro,
amino that may be substituted, imino that may be
20 substituted, cyano, mercapto, monosubstituted thio that
may be substituted, monosubstituted thioxy that may be
substituted, monosubstituted sulfinyl that may be
substituted, monosubstituted sulfonyl that may be
substituted, sulfo, or trisubstituted silyl, and any
25 combinations of R_{1a} , R_{2a} , R_3-R_5 may together form a ring
structure.

20. A NAD(P)H oxydase inhibitor, comprising
a compound represented by the formula (I):



or a prodrug, pharmaceutically acceptable salt or
5 solvate thereof,

wherein, in the formula,

R₁ is hydrogen, lower alkyl, amino that
may be substituted, or aryl lower alkyl that may be
substituted; and

10 R₂ is hydrogen, lower alkyl that may be
substituted, cycloalkyl that may be substituted,
cycloalkyl lower alkyl that may be substituted, lower
alkoxy that may be substituted, aryl that may be
substituted, aryl lower alkyl that may be substituted,
15 aryloxy lower alkyl that may be substituted, lower
alkylsulfonyl that may be substituted, arylsulfonyl
that may be substituted, heteroaryl lower alkyl that
may be substituted, heterocyclic group lower alkyl that
may be substituted, or amino that may be substituted;
20 or

R₁ and R₂ together with adjacent N atom
may form a heterocycle that may be substituted;

R₃ is hydrogen, hydroxy, lower alkoxy,
halogen, or amino that may be substituted;

25 R₄ is hydrogen, lower alkyl, or aryl that
may be substituted; and

R₅ is hydroxy, lower alkyl that may be substituted, aryl that may be substituted, aryl lower alkyl that may be substituted, cycloalkyl lower alkyl
30 that may be substituted, aryl lower alkenyl that may be substituted, cycloalkyl lower alkenyl that may be substituted, aryl lower alkynyl that may be substituted, cycloalkyl lower alkynyl that may be substituted, aryl carbonyl that may be substituted,
35 aryl lower alkyl carbonyl that may be substituted, heterocyclic group that may be substituted, halogen, CHO, amino that may be substituted, or imino that may be substituted.

21. A method of preventing or treating NAD(P)H-related diseases, comprising administering the compound of any one of claims 1-20 to an animal including human.

22. The method of claim 21, wherein said disease is selected from the group consisting of inflammation, pulmonary circulation disorders, ischemic heart disease, cerebral circulation disorders,
5 arteriosclerosis, diabetic complications, hypertension, and proliferative disorders.

23. The method of claim 21, wherein said disease is brain infarction or diabetic retinal disorder.

24. A use of the compound of any one of claims 1-20 for the manufacture of pharmaceuticals employed for preventing or treating NAD(P)H-related diseases.

25. The use of claim 24, wherein said disease is selected from the group consisting of inflammation, pulmonary circulation disorders, ischemic heart disease, cerebral circulation disorders,
5 arteriosclerosis, diabetic complications, hypertension, and proliferative disorders.

26. The use of claim 24, wherein said disease is brain infarction or diabetic retinal disorder.

27. The compound of claim 1, or a prodrug, pharmaceutically acceptable salt or solvate thereof, wherein R_{1a} is carbamoyl that may be substituted.

28. The compound of claim 1, or a prodrug, pharmaceutically acceptable salt or solvate thereof, wherein R_{1a} is carbamoyl that may be substituted, and R_{2a} is hydrogen.

29. A medicament, comprising the compound of claims 27 or 28.

30. A NAD(P)H oxydase inhibitor, comprising the compound of claims 27 or 28.